

Acta Crystallographica Section E

Structure Reports

Online

ISSN 1600-5368

trans,trans,trans-Diaquabis(nicotinamide- κ N)bis(2-nitrobenzoato- κ O)cadmium(II) dihydrate

Kou-Lin Zhang,^a Bo Yang,^a Jian-Guo Lin^a and Seik Weng Ng^{b*}

^aCollege of Chemistry and Chemical Engineering, Yangzhou University, Yangzhou 225002, People's Republic of China, and ^bDepartment of Chemistry, University of Malaya, 50603, Kuala Lumpur, Malaysia
Correspondence e-mail: seikweng@um.edu.my

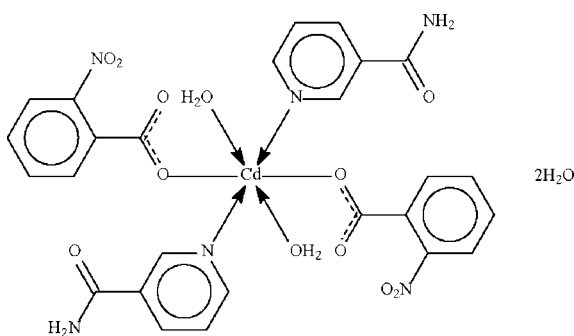
Received 21 January 2009; accepted 16 February 2009

Key indicators: single-crystal X-ray study; $T = 293$ K; mean $\sigma(\text{C}-\text{C}) = 0.005$ Å; R factor = 0.034; wR factor = 0.093; data-to-parameter ratio = 10.8.

The cadmium atom in the title compound, $[\text{Cd}(\text{C}_7\text{H}_4\text{NO}_4)_2(\text{C}_6\text{H}_6\text{N}_2\text{O})_2(\text{H}_2\text{O})_2] \cdot 2\text{H}_2\text{O}$, lies on a center of inversion in an all-*trans* octahedral environment. In the crystal, the complex interacts with the uncoordinated water molecules through O—H...O and N—H...O hydrogen bonds, forming a layered network.

Related literature

There are several examples of diaquadi(arylcarboxylato)di(nicotinamide)metal(II) compounds. For recent examples, see: Hökelek & Necefoğlu (2007*a,b*); Hökelek *et al.* (2007); Koksharova *et al.* (2006); Şahin *et al.* (2007*a,b*); Stachova *et al.* (2006); Çaylak *et al.* (2007).



Experimental

Crystal data

$[\text{Cd}(\text{C}_7\text{H}_4\text{NO}_4)_2(\text{C}_6\text{H}_6\text{N}_2\text{O})_2(\text{H}_2\text{O})_2] \cdot 2\text{H}_2\text{O}$
 $M_r = 760.94$
Monoclinic, $P2_1/n$
 $a = 7.9365$ (8) Å

$b = 19.589$ (2) Å
 $c = 10.059$ (1) Å
 $\beta = 103.178$ (2)°
 $V = 1522.6$ (3) Å³
 $Z = 2$

Mo $K\alpha$ radiation
 $\mu = 0.80$ mm⁻¹

$T = 293$ K
 $0.50 \times 0.18 \times 0.18$ mm

Data collection

Bruker SMART area-detector diffractometer
Absorption correction: multi-scan (SADABS; Sheldrick, 1996)
 $T_{\min} = 0.619$, $T_{\max} = 0.866$

4427 measured reflections
2651 independent reflections
2396 reflections with $I > 2\sigma(I)$
 $R_{\text{int}} = 0.016$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.034$
 $wR(F^2) = 0.093$
 $S = 1.13$
2651 reflections
246 parameters
9 restraints

H atoms treated by a mixture of independent and constrained refinement
 $\Delta\rho_{\max} = 0.42$ e Å⁻³
 $\Delta\rho_{\min} = -1.04$ e Å⁻³

Table 1

Hydrogen-bond geometry (Å, °).

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
O1w—H11...O2w	0.85 (4)	1.92 (4)	2.764 (4)	174 (4)
O1w—H12...O2 ⁱ	0.85 (4)	1.95 (5)	2.718 (4)	150 (4)
O2w—H21...O5 ⁱ	0.85 (3)	2.08 (3)	2.910 (4)	166 (4)
O2w—H22...O1 ⁱⁱ	0.85 (3)	2.00 (1)	2.846 (3)	177 (5)
N3—H31...O2 ⁱⁱⁱ	0.85 (3)	2.22 (2)	3.038 (4)	165 (4)
N3—H32...O5 ^{iv}	0.85 (3)	2.05 (3)	2.873 (4)	164 (4)

Symmetry codes: (i) $-x+1, -y+1, -z+1$; (ii) $-x+2, -y+1, -z+1$; (iii) $x, y, z+1$; (iv) $-x, -y+1, -z+2$.

Data collection: SMART (Bruker, 2000); cell refinement: SAINT (Bruker, 2000); data reduction: SAINT; program(s) used to solve structure: SHELXS97 (Sheldrick, 2008); program(s) used to refine structure: SHELXL97 (Sheldrick, 2008); molecular graphics: X-SEED (Barbour, 2001); software used to prepare material for publication: publCIF (Westrip, 2009).

We thank the Foundation of Jiangsu Provincial Key Program of Physical Chemistry in Yangzhou University and the University of Malaya for supporting this study.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: PV2136).

References

- Barbour, L. J. (2001). *J. Supramol. Chem.*, **1**, 189–191.
Bruker (2000). SAINT and S.A.I.N.T. Bruker AXS Inc., Madison, Wisconsin, USA.
Çaylak, N., Hökelek, T. & Necefoğlu, H. (2007). *Acta Cryst.* **E63**, m1341–m1343.
Hökelek, T., Çaylak, N. & Necefoğlu, H. (2007). *Acta Cryst.* **E63**, m1873–m1874.
Hökelek, T. & Necefoğlu, H. (2007*a*). *Acta Cryst.* **E63**, m1078–m1080.
Hökelek, T. & Necefoğlu, H. (2007*b*). *Acta Cryst.* **E63**, m1279–m1281.
Koksharova, T. V., Sadikov, G. G., Antsyshkina, A. S., Gritsenko, I. S., Sergienko, V. S. & Egorova, O. A. (2006). *Russ. J. Inorg. Chem.* **51**, 895–900.
Şahin, O., Büyükgüngör, O., Köse, D. A. & Necefoglu, H. (2007*a*). *Acta Cryst.* **C63**, m510–m512.
Şahin, O., Büyükgüngör, O., Köse, D. A., Ozturkkan, E. F. & Necefoglu, H. (2007*b*). *Acta Cryst.* **C63**, m243–m245.
Sheldrick, G. M. (1996). SADABS. University of Göttingen, Germany.
Sheldrick, G. M. (2008). *Acta Cryst.* **A64**, 112–122.
Stachova, P., Melnik, M., Korobik, M., Mrozinski, M., Koman, M., Glowiak, T. & Valigura, D. (2006). *Inorg. Chim. Acta*, **360**, 1517–1522.
Westrip, S. P. (2009). publCIF. In preparation.

supporting information

Acta Cryst. (2009). E65, m292 [doi:10.1107/S1600536809005479]

***trans,trans,trans*-Diaquabis(nicotinamide- κ N)bis(2-nitrobenzoato- κ O)cadmium(II) dihydrate**

Kou-Lin Zhang, Bo Yang, Jian-Guo Lin and Seik Weng Ng

S1. Experimental

A water/methanol (1:1 v/v) solution (3 ml) of cadmium nitrate trihydrate (0.082 g, 0.3 mmol) was added to a water/methanol (1:1 v/v) solution (3 ml) of 2-nitrobenzoic acid (0.100 g, 0.6 mmol), sodium hydroxide (0.024 g, 0.6 mmol) and nicotinamide (0.073 g, 0.6 mmol). A white powder was obtained after several days; this was recrystallized from DMF/methanol (3:1 v/v) to give colorless crystals in 50% yield. CH&N elemental analysis. Calculated for $C_{26}H_{28}CdN_6O_{14}$: C 41.04 H 3.68 N 11.04%; found: C 40.08, H 3.87, N 10.93%.

S2. Refinement

Carbon-bound H atoms were placed in calculated positions and were allowed to ride on the parent atoms. N and O-bound H atoms were located in a difference Fourier map, and were refined with distance restraints $N-H = O-H = 0.85 \pm 0.01$ Å; for the water molecules, an additional $H \cdots H$ 1.39 ± 0.01 Å restraint was used. Their temperature factors were freely refined.

The measurements are 100% at the 2θ limit of 50° .

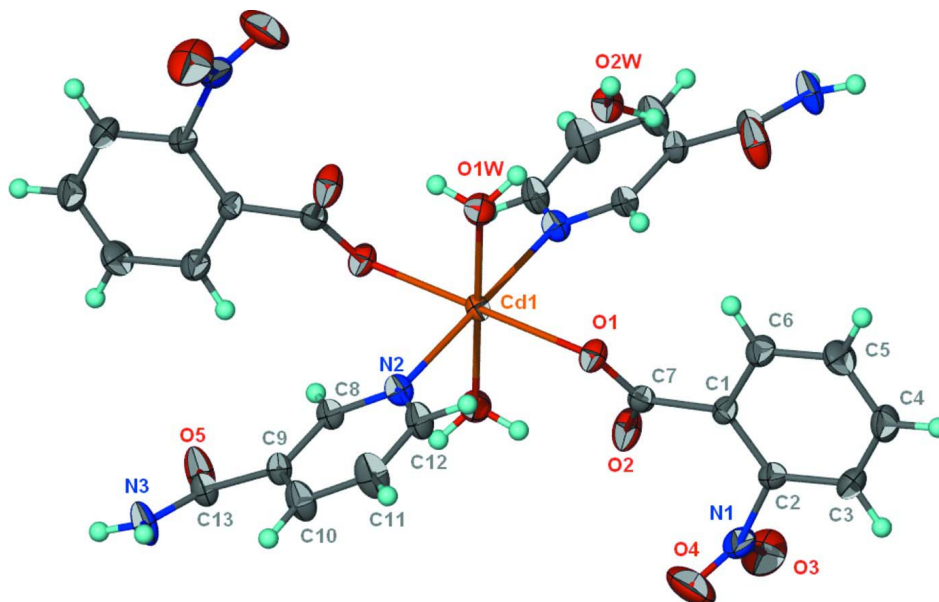


Figure 1

Thermal ellipsoid plot of $Cd(H_2O)_2(C_7H_4NO_4)_2(C_6H_6N_2O)_2 \cdot 2H_2O$; displacement ellipsoids are drawn at the 50% probability level, and H atoms as spheres of arbitrary radii.

***trans,trans,trans*-Diaquabis(nicotinamide- κ N)bis(2- nitrobenzoato- κ O)cadmium(II) dihydrate**

Crystal data

[Cd(C₇H₄NO₄)₂(C₆H₆N₂O)₂(H₂O)₂] \cdot 2H₂O
M_r = 760.94
 Monoclinic, *P*2₁/*n*
 Hall symbol: -*P* 2yn
a = 7.9365 (8) Å
b = 19.589 (2) Å
c = 10.059 (1) Å
 β = 103.178 (2)°
V = 1522.6 (3) Å³
Z = 2

F(000) = 772
D_x = 1.660 Mg m⁻³
 Mo *K* α radiation, λ = 0.71073 Å
 Cell parameters from 3417 reflections
 θ = 2.1–25.1°
 μ = 0.80 mm⁻¹
T = 293 K
 Rod, colorless
 0.50 \times 0.18 \times 0.18 mm

Data collection

Bruker SMART area-detector
 diffractometer
 Radiation source: medium-focus sealed tube
 Graphite monochromator
 φ and ω scans
 Absorption correction: multi-scan
 (*SADABS*; Sheldrick, 1996)
T_{min} = 0.619, *T_{max}* = 0.866

4427 measured reflections
 2651 independent reflections
 2396 reflections with *I* > 2 σ (*I*)
R_{int} = 0.016
 θ_{\max} = 25.1°, θ_{\min} = 2.1°
h = -9→3
k = -20→23
l = -11→11

Refinement

Refinement on *F*²
 Least-squares matrix: full
R[*F*² > 2 σ (*F*²)] = 0.034
wR(*F*²) = 0.093
S = 1.13
 2651 reflections
 246 parameters
 9 restraints
 Primary atom site location: structure-invariant
 direct methods

Secondary atom site location: difference Fourier
 map
 Hydrogen site location: inferred from
 neighbouring sites
 H atoms treated by a mixture of independent
 and constrained refinement
 $w = 1/[\sigma^2(F_o^2) + (0.0433P)^2 + 2.3534P]$
 where $P = (F_o^2 + 2F_c^2)/3$
 $(\Delta/\sigma)_{\max} = 0.001$
 $\Delta\rho_{\max} = 0.42 \text{ e } \text{Å}^{-3}$
 $\Delta\rho_{\min} = -1.04 \text{ e } \text{Å}^{-3}$

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters (Å²)

	<i>x</i>	<i>y</i>	<i>z</i>	<i>U_{iso}</i> */ <i>U_{eq}</i>
Cd1	0.5000	0.5000	0.5000	0.02349 (13)
O1	0.6570 (3)	0.40982 (12)	0.4385 (2)	0.0312 (5)
O2	0.4256 (3)	0.36571 (14)	0.2970 (3)	0.0428 (7)
O3	0.3595 (4)	0.23243 (19)	0.1217 (4)	0.0693 (10)
O4	0.4552 (5)	0.2124 (2)	0.3355 (4)	0.0835 (12)
O5	0.0789 (4)	0.48477 (17)	0.8538 (3)	0.0505 (8)
O1W	0.7685 (3)	0.55083 (13)	0.5819 (3)	0.0362 (6)
H11	0.842 (5)	0.561 (2)	0.536 (4)	0.055 (14)*
H12	0.737 (6)	0.5868 (14)	0.616 (5)	0.082 (19)*
O2W	0.9923 (3)	0.57871 (14)	0.4149 (3)	0.0375 (6)
H21	0.987 (5)	0.556 (2)	0.342 (3)	0.079 (18)*
H22	1.097 (2)	0.581 (2)	0.460 (3)	0.050 (13)*

N1	0.4713 (4)	0.23414 (16)	0.2260 (4)	0.0438 (8)
N2	0.5129 (3)	0.44979 (15)	0.7119 (3)	0.0291 (6)
N3	0.1940 (4)	0.44256 (19)	1.0627 (3)	0.0405 (8)
H31	0.274 (4)	0.424 (2)	1.122 (3)	0.055 (13)*
H32	0.114 (4)	0.457 (2)	1.098 (3)	0.049 (12)*
C1	0.6994 (4)	0.32348 (16)	0.2832 (3)	0.0237 (6)
C2	0.6421 (4)	0.26190 (17)	0.2198 (3)	0.0282 (7)
C3	0.7386 (5)	0.22316 (19)	0.1503 (4)	0.0386 (9)
H3	0.6942	0.1829	0.1070	0.047 (12)*
C4	0.9033 (5)	0.2454 (2)	0.1459 (4)	0.0407 (9)
H4	0.9714	0.2196	0.1008	0.047 (12)*
C5	0.9655 (5)	0.3057 (2)	0.2087 (4)	0.0430 (9)
H5	1.0761	0.3205	0.2063	0.057 (13)*
C6	0.8643 (4)	0.34455 (19)	0.2755 (4)	0.0337 (8)
H6	0.9075	0.3855	0.3161	0.046 (12)*
C7	0.5842 (4)	0.36885 (16)	0.3458 (3)	0.0262 (7)
C8	0.3792 (4)	0.46351 (17)	0.7674 (3)	0.0256 (7)
H8	0.2975	0.4951	0.7238	0.044 (12)*
C9	0.3553 (4)	0.43356 (17)	0.8857 (3)	0.0273 (7)
C10	0.4767 (5)	0.3864 (2)	0.9499 (4)	0.0399 (9)
H10	0.4645	0.3648	1.0295	0.044 (11)*
C11	0.6165 (5)	0.3721 (2)	0.8941 (4)	0.0461 (10)
H11A	0.7004	0.3410	0.9362	0.055 (13)*
C12	0.6301 (5)	0.40434 (19)	0.7751 (4)	0.0352 (8)
H12A	0.7239	0.3942	0.7375	0.052 (12)*
C13	0.1972 (4)	0.45472 (18)	0.9337 (3)	0.0312 (7)

Atomic displacement parameters (\AA^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Cd1	0.02157 (19)	0.0277 (2)	0.0224 (2)	-0.00033 (12)	0.00743 (13)	0.00046 (12)
O1	0.0292 (12)	0.0324 (13)	0.0303 (12)	0.0020 (10)	0.0034 (10)	-0.0075 (10)
O2	0.0268 (13)	0.0550 (17)	0.0443 (15)	0.0049 (12)	0.0033 (11)	-0.0207 (13)
O3	0.0398 (17)	0.078 (2)	0.085 (3)	-0.0141 (16)	0.0042 (17)	-0.027 (2)
O4	0.087 (3)	0.098 (3)	0.078 (3)	-0.041 (2)	0.044 (2)	0.003 (2)
O5	0.0388 (16)	0.084 (2)	0.0318 (15)	0.0297 (15)	0.0135 (12)	0.0122 (14)
O1W	0.0273 (13)	0.0406 (14)	0.0416 (15)	-0.0066 (11)	0.0094 (11)	-0.0085 (12)
O2W	0.0276 (13)	0.0452 (15)	0.0390 (15)	0.0019 (11)	0.0061 (11)	-0.0039 (12)
N1	0.0417 (19)	0.0363 (18)	0.057 (2)	-0.0114 (14)	0.0190 (17)	-0.0155 (16)
N2	0.0251 (14)	0.0348 (15)	0.0274 (15)	0.0037 (12)	0.0059 (11)	0.0027 (12)
N3	0.0366 (17)	0.066 (2)	0.0223 (15)	0.0158 (16)	0.0128 (13)	0.0093 (15)
C1	0.0281 (16)	0.0240 (16)	0.0173 (14)	0.0011 (13)	0.0015 (12)	-0.0017 (12)
C2	0.0282 (17)	0.0258 (17)	0.0313 (17)	-0.0024 (13)	0.0083 (14)	-0.0016 (14)
C3	0.051 (2)	0.0263 (18)	0.041 (2)	-0.0017 (16)	0.0154 (17)	-0.0088 (16)
C4	0.042 (2)	0.040 (2)	0.045 (2)	0.0102 (17)	0.0185 (17)	-0.0024 (17)
C5	0.0305 (19)	0.047 (2)	0.056 (3)	-0.0013 (17)	0.0203 (18)	-0.0033 (19)
C6	0.0303 (18)	0.0349 (19)	0.0359 (19)	-0.0044 (15)	0.0075 (15)	-0.0083 (15)
C7	0.0254 (16)	0.0243 (16)	0.0297 (17)	0.0019 (13)	0.0078 (13)	0.0013 (13)

C8	0.0246 (16)	0.0309 (18)	0.0201 (15)	0.0033 (13)	0.0025 (12)	0.0029 (13)
C9	0.0282 (17)	0.0336 (18)	0.0199 (15)	0.0028 (14)	0.0051 (13)	-0.0012 (13)
C10	0.042 (2)	0.053 (2)	0.0267 (18)	0.0173 (18)	0.0117 (15)	0.0150 (17)
C11	0.041 (2)	0.062 (3)	0.035 (2)	0.027 (2)	0.0095 (17)	0.0190 (19)
C12	0.0303 (18)	0.046 (2)	0.0317 (19)	0.0088 (16)	0.0117 (15)	0.0022 (16)
C13	0.0304 (18)	0.0375 (19)	0.0274 (17)	0.0051 (15)	0.0098 (14)	0.0027 (14)

Geometric parameters (Å, °)

Cd1—O1 ⁱ	2.325 (2)	C1—C2	1.391 (4)
Cd1—O1	2.325 (2)	C1—C6	1.391 (5)
Cd1—O1W ⁱ	2.326 (2)	C1—C7	1.512 (4)
Cd1—O1W	2.326 (2)	C2—C3	1.377 (5)
Cd1—N2	2.329 (3)	C3—C4	1.388 (5)
Cd1—N2 ⁱ	2.329 (3)	C3—H3	0.9300
O1—C7	1.266 (4)	C4—C5	1.378 (6)
O2—C7	1.244 (4)	C4—H4	0.9300
O3—N1	1.211 (5)	C5—C6	1.386 (5)
O4—N1	1.214 (5)	C5—H5	0.9300
O5—C13	1.237 (4)	C6—H6	0.9300
O1W—H11	0.85 (4)	C8—C9	1.378 (5)
O1W—H12	0.85 (4)	C8—H8	0.9300
O2W—H21	0.85 (3)	C9—C10	1.383 (5)
O2W—H22	0.85 (3)	C9—C13	1.502 (5)
N1—C2	1.476 (4)	C10—C11	1.382 (5)
N2—C8	1.334 (4)	C10—H10	0.9300
N2—C12	1.339 (4)	C11—C12	1.379 (5)
N3—C13	1.325 (4)	C11—H11A	0.9300
N3—H31	0.84 (3)	C12—H12A	0.9300
N3—H32	0.85 (3)		
O1 ⁱ —Cd1—O1	180.000 (1)	C1—C2—N1	120.5 (3)
O1 ⁱ —Cd1—O1W ⁱ	85.20 (9)	C2—C3—C4	118.7 (3)
O1—Cd1—O1W ⁱ	94.80 (9)	C2—C3—H3	120.7
O1 ⁱ —Cd1—O1W	94.80 (9)	C4—C3—H3	120.7
O1—Cd1—O1W	85.20 (9)	C5—C4—C3	119.8 (3)
O1W ⁱ —Cd1—O1W	180.00 (12)	C5—C4—H4	120.1
O1 ⁱ —Cd1—N2	89.52 (9)	C3—C4—H4	120.1
O1—Cd1—N2	90.48 (9)	C4—C5—C6	120.4 (3)
O1W ⁱ —Cd1—N2	89.36 (10)	C4—C5—H5	119.8
O1W—Cd1—N2	90.64 (10)	C6—C5—H5	119.8
O1 ⁱ —Cd1—N2 ⁱ	90.48 (9)	C5—C6—C1	121.4 (3)
O1—Cd1—N2 ⁱ	89.52 (9)	C5—C6—H6	119.3
O1W ⁱ —Cd1—N2 ⁱ	90.64 (10)	C1—C6—H6	119.3
O1W—Cd1—N2 ⁱ	89.36 (10)	O2—C7—O1	125.0 (3)
N2—Cd1—N2 ⁱ	180.0	O2—C7—C1	117.4 (3)
C7—O1—Cd1	119.7 (2)	O1—C7—C1	117.5 (3)
Cd1—O1W—H11	126 (3)	N2—C8—C9	123.7 (3)

Cd1—O1W—H12	100 (3)	N2—C8—H8	118.2
H11—O1W—H12	109 (4)	C9—C8—H8	118.2
H21—O2W—H22	109.4 (17)	C8—C9—C10	118.0 (3)
O3—N1—O4	124.7 (4)	C8—C9—C13	116.7 (3)
O3—N1—C2	118.2 (4)	C10—C9—C13	125.3 (3)
O4—N1—C2	117.1 (4)	C11—C10—C9	118.9 (3)
C8—N2—C12	117.9 (3)	C11—C10—H10	120.5
C8—N2—Cd1	115.1 (2)	C9—C10—H10	120.5
C12—N2—Cd1	126.6 (2)	C12—C11—C10	119.3 (3)
C13—N3—H31	126 (3)	C12—C11—H11A	120.4
C13—N3—H32	122 (2)	C10—C11—H11A	120.4
H31—N3—H32	111.4 (18)	N2—C12—C11	122.2 (3)
C2—C1—C6	116.4 (3)	N2—C12—H12A	118.9
C2—C1—C7	122.4 (3)	C11—C12—H12A	118.9
C6—C1—C7	121.0 (3)	O5—C13—N3	122.8 (3)
C3—C2—C1	123.3 (3)	O5—C13—C9	119.2 (3)
C3—C2—N1	116.2 (3)	N3—C13—C9	118.0 (3)
O1W ⁱ —Cd1—O1—C7	22.2 (2)	C4—C5—C6—C1	1.0 (6)
O1W—Cd1—O1—C7	-157.8 (2)	C2—C1—C6—C5	-0.3 (5)
N2—Cd1—O1—C7	111.6 (2)	C7—C1—C6—C5	-174.8 (3)
N2 ⁱ —Cd1—O1—C7	-68.4 (2)	Cd1—O1—C7—O2	-11.8 (5)
O1 ⁱ —Cd1—N2—C8	29.7 (2)	Cd1—O1—C7—C1	163.9 (2)
O1—Cd1—N2—C8	-150.3 (2)	C2—C1—C7—O2	-27.5 (5)
O1W ⁱ —Cd1—N2—C8	-55.5 (2)	C6—C1—C7—O2	146.7 (3)
O1W—Cd1—N2—C8	124.5 (2)	C2—C1—C7—O1	156.4 (3)
O1 ⁱ —Cd1—N2—C12	-157.2 (3)	C6—C1—C7—O1	-29.4 (4)
O1—Cd1—N2—C12	22.8 (3)	C12—N2—C8—C9	0.0 (5)
O1W ⁱ —Cd1—N2—C12	117.6 (3)	Cd1—N2—C8—C9	173.7 (3)
O1W—Cd1—N2—C12	-62.4 (3)	N2—C8—C9—C10	-0.3 (5)
C6—C1—C2—C3	-1.2 (5)	N2—C8—C9—C13	-179.8 (3)
C7—C1—C2—C3	173.2 (3)	C8—C9—C10—C11	0.7 (6)
C6—C1—C2—N1	178.0 (3)	C13—C9—C10—C11	-179.9 (4)
C7—C1—C2—N1	-7.5 (5)	C9—C10—C11—C12	-0.8 (7)
O3—N1—C2—C3	-69.1 (5)	C8—N2—C12—C11	-0.1 (6)
O4—N1—C2—C3	108.4 (4)	Cd1—N2—C12—C11	-173.0 (3)
O3—N1—C2—C1	111.6 (4)	C10—C11—C12—N2	0.6 (7)
O4—N1—C2—C1	-70.9 (5)	C8—C9—C13—O5	15.8 (5)
C1—C2—C3—C4	1.9 (6)	C10—C9—C13—O5	-163.7 (4)
N1—C2—C3—C4	-177.4 (3)	C8—C9—C13—N3	-161.5 (3)
C2—C3—C4—C5	-1.1 (6)	C10—C9—C13—N3	19.0 (6)
C3—C4—C5—C6	-0.3 (6)		

Symmetry code: (i) $-x+1, -y+1, -z+1$.

Hydrogen-bond geometry (\AA , $^\circ$)

$D-H\cdots A$	$D-H$	$H\cdots A$	$D\cdots A$	$D-H\cdots A$
O1w—H11 \cdots O2w	0.85 (4)	1.92 (4)	2.764 (4)	174 (4)

O1 _w —H12...O2 ⁱ	0.85 (4)	1.95 (5)	2.718 (4)	150 (4)
O2 _w —H21...O5 ⁱ	0.85 (3)	2.08 (3)	2.910 (4)	166 (4)
O2 _w —H22...O1 ⁱⁱ	0.85 (3)	2.00 (1)	2.846 (3)	177 (5)
N3—H31...O2 ⁱⁱⁱ	0.85 (3)	2.22 (2)	3.038 (4)	165 (4)
N3—H32...O5 ^{iv}	0.85 (3)	2.05 (3)	2.873 (4)	164 (4)

Symmetry codes: (i) $-x+1, -y+1, -z+1$; (ii) $-x+2, -y+1, -z+1$; (iii) $x, y, z+1$; (iv) $-x, -y+1, -z+2$.